Trying 3106016892...Open

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* The files listed above are temporarily unavailable.

FILE 'HOME' ENTERED AT 15:30:49 ON 16 AUG 2000

09/270,006

=> file reg

COST IN U.S. DOLLARS

SINCE FILE

TOTAL

FULL ESTIMATED COST

ENTRY 0.15

SESSION 0.15

FILE 'REGISTRY' ENTERED AT 15:30:55 ON 16 AUG 2000 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS. COPYRIGHT (C) 2000 American Chemical Society (ACS)

STRUCTURE FILE UPDATES: 15 AUG 2000 HIGHEST RN 286364-83-8 DICTIONARY FILE UPDATES: 15 AUG 2000 HIGHEST RN 286364-83-8

TSCA INFORMATION NOW CURRENT THROUGH JANUARY 11, 2000

Please note that search-term pricing does apply when conducting SmartSELECT searches.

Structure search limits have been increased. See HELP SLIMIT for details.

=>Testing the current file.... screen

ENTER SCREEN EXPRESSION OR (END):end

=> screen 2076

L1 SCREEN CREATED

=>

Uploading c:\stnexp4\queries\09270006.str

L2 STRUCTURE UPLOADED

=> que L2 AND L1

L3 QUE L2 AND L1

=> d 12

L2 HAS NO ANSWERS

L2

STR



Structure attributes must be viewed using STN Express query preparation.

=> s 12 sss sam

SAMPLE SEARCH INITIATED 15:31:35 FILE 'REGISTRY' SAMPLE SCREEN SEARCH COMPLETED - 47148 TO ITERATE

2.1% PROCESSED 1000 ITERATIONS INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED) SEARCH TIME: 00.00.01

3 ANSWERS

FULL FILE PROJECTIONS: ONLINE **INCOMPLETE**

BATCH **INCOMPLETE**

PROJECTED ITERATIONS:

930168 TO 955752

PROJECTED ANSWERS:

2115 TO 3541

L4

3 SEA SSS SAM L2

=> d sacn

'SACN' IS NOT A VALID FORMAT FOR FILE 'REGISTRY'

The following are valid formats:

Substance information can be displayed by requesting individual fields or predefined formats. The predefined substance formats are: (RN = CAS Registry Number)

REG - RN

SAM - Index Name, MF, and structure - no RN FIDE - All substance data, except sequence data

IDE - FIDE, but only 50 names SQIDE - IDE, plus sequence data

SQIDE3 - Same as SQIDE, but 3-letter amino acid codes are used

SQD - Protein sequence data, includes RN

SQD3 - Same as SQD, but 3-letter amino acid codes are used

SQN - Protein sequence name information, includes RN

Any CA File format may be combined with any substance format to obtain CA references citing the substance. The substance formats must be cited first. The CA File predefined formats are:

ABS -- Abstract

APPS -- Application and Priority Information

BIB -- CA Accession Number, plus Bibliographic Data

CAN -- CA Accession Number

CBIB -- CA Accession Number, plus Bibliographic Data (compressed)

IND -- Index Data

IPC -- International Patent Classification

PATS -- PI, SO

STD -- BIB, IPC, and NCL

IABS --ABS, indented, with text labels

IBIB -- BIB, indented, with text labels

ISTD -- STD format, indented

OBIB ----- AN, plus Bibliographic Data (original)

OIBIB ---- OBIB, indented with text labels

SBIB ----- BIB, no citations SIBIB ----- IBIB, no citations

09/270,006

The ALL format gives FIDE BIB ABS IND RE, plus sequence data when it is available.

The MAX format is the same as ALL.

The IALL format is the same as ALL with BIB ABS and IND indented, with text labels.

For additional information, please consult the following help messages:

HELP DFIELDS -- To see a complete list of individual display fields. HELP FORMATS -- To see detailed descriptions of the predefined formats. ENTER DISPLAY FORMAT (IDE):d scan

'D' IS NOT A VALID FORMAT FOR FILE 'REGISTRY' 'SCAN' IS NOT A VALID FORMAT FOR FILE 'REGISTRY'

The following are valid formats:

Substance information can be displayed by requesting individual fields or predefined formats. The predefined substance formats are: (RN = CAS Registry Number)

REG - RN

SAM - Index Name, MF, and structure - no RN
FIDE - All substance data, except sequence data

IDE - FIDE, but only 50 names SQIDE - IDE, plus sequence data

SQIDE3 - Same as SQIDE, but 3-letter amino acid codes are used

SQD - Protein sequence data, includes RN

SQD3 - Same as SQD, but 3-letter amino acid codes are used

SQN - Protein sequence name information, includes RN

Any CA File format may be combined with any substance format to obtain CA references citing the substance. The substance formats must be cited first. The CA File predefined formats are:

ABS -- Abstract

APPS -- Application and Priority Information

BIB -- CA Accession Number, plus Bibliographic Data

CAN -- CA Accession Number

CBIB -- CA Accession Number, plus Bibliographic Data (compressed)

IND -- Index Data

IPC -- International Patent Classification

PATS -- PI, SO

STD -- BIB, IPC, and NCL

IABS --ABS, indented, with text labels

IBIB -- BIB, indented, with text labels

ISTD -- STD format, indented

OBIB ----- AN, plus Bibliographic Data (original)

OIBIB ----- OBIB, indented with text labels

SBIB ----- BIB, no citations SIBIB ----- IBIB, no citations

The ALL format gives FIDE BIB ABS IND RE, plus sequence data when it is available.

The MAX format is the same as ALL.

The IALL format is the same as ALL with BIB ABS and IND indented, with text labels.

For additional information, please consult the following help messages:

HELP DFIELDS -- To see a complete list of individual display fields. HELP FORMATS -- To see detailed descriptions of the predefined formats. ENTER DISPLAY FORMAT (IDE):end

=> d scan

L4 3 ANSWERS REGISTRY COPYRIGHT 2000 ACS

IN 2-Furancarboxylic acid, 2-[4-[[[[1-[(ethoxyhydroxyphosphino)methyl]-1H-

MF C14 H19 N6 O9 P S . 2 Na

Absolute stereochemistry.

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):2

L4 3 ANSWERS REGISTRY COPYRIGHT 2000 ACS
IN Oxiranecarboxylic acid, 3-(4-chlorophenyl)-, sodium salt, trans- (9CI)
09/270,006 Page 5

MF C9 H7 Cl O3 . Na

Relative stereochemistry.

Na

L43 ANSWERS REGISTRY COPYRIGHT 2000 ACS

IN Oxiranecarboxylic acid, 3-[[[(1S)-2-oxo-2-[(2-phenylethyl)amino]-1-(phenylmethyl)ethyl]amino]carbonyl]-, (2S,3S)- (9CI)

C21 H22 N2 O5 MF

Absolute stereochemistry.

ALL ANSWERS HAVE BEEN SCANNED

 \Rightarrow log y

COST IN U.S. DOLLARS

SINCE FILE TOTAL ENTRY SESSION 0.90 1.05

FULL ESTIMATED COST

STN INTERNATIONAL LOGOFF AT 15:32:33 ON 16 AUG 2000

Trying 3106016892...Open

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| | | | | BY ADDITION OF CITED REFERENCES TO CAPLUS, CA, |
| | | | | REGISTRY, CASREACT, MARPAT, and MARPATPREV |
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| NEWS | 14 | Jul | 27 | Derwent Journal Of Synthetic Methods Reloaded |
| | | | | with New Data |
| NEWS | 15 | Jul | 27 | DERWENT WORLD PATENTS INDEX: FAST TRACK RELEASE OF |
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* The files listed above are temporarily unavailable.

FILE 'HOME' ENTERED AT 15:37:06 ON 16 AUG 2000

=> file reg

COST IN U.S. DOLLARS
SINCE FILE TOTAL
ENTRY SESSION
FULL ESTIMATED COST
0.15

FILE 'REGISTRY' ENTERED AT 15:37:12 ON 16 AUG 2000 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. 09/270,006

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STRUCTURE FILE UPDATES: 15 AUG 2000 HIGHEST RN 286364-83-8 DICTIONARY FILE UPDATES: 15 AUG 2000 HIGHEST RN 286364-83-8

TSCA INFORMATION NOW CURRENT THROUGH JANUARY 11, 2000

Please note that search-term pricing does apply when conducting SmartSELECT searches.

Structure search limits have been increased. See HELP SLIMIT for details.

=>Testing the current file.... screen

ENTER SCREEN EXPRESSION OR (END):end

=> screen 2076

L1 SCREEN CREATED

=>

Uploading c:\stnexp4\queries\09270006.str

L2 STRUCTURE UPLOADED

=> que L2 AND L1

L3 QUE L2 AND L1

=> d 12

L2 HAS NO ANSWERS

L2 STR

$$G2$$
 $G2$
 $G1$
 CH_2
 $O-8$
 $G2$

G1 O, S, P

G2 H, Cy, Hy, Ak

Structure attributes must be viewed using STN Express query preparation.

=> s 12 sss sam

SAMPLE SEARCH INITIATED 15:37:52 FILE 'REGISTRY' 09/270,006

SAMPLE SCREEN SEARCH COMPLETED - 47148 TO ITERATE

2.1% PROCESSED 1000 ITERATIONS

INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **INCOMPLETE**

BATCH **INCOMPLETE**

PROJECTED ITERATIONS:

930168 TO 955752

PROJECTED ANSWERS:

1303 TO 2467

L4

2 SEA SSS SAM L2

=> d scan

L4 2 ANSWERS REGISTRY COPYRIGHT 2000 ACS

IN Oxiranecarboxylic acid, 3-(4-chlorophenyl)-, sodium salt, trans- (9CI)

MF C9 H7 C1 O3 . Na

Relative stereochemistry.

Na

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L4 2 ANSWERS REGISTRY COPYRIGHT 2000 ACS

IN Oxiranecarboxylic acid, 3-[[[(1S)-2-oxo-2-[(2-phenylethyl)amino]-1-

(phenylmethyl)ethyl]amino]carbonyl]-, (2S,3S)- (9CI)

MF C21 H22 N2 O5

Absolute stereochemistry.

09/270,006

2 ANSWERS

ALL ANSWERS HAVE BEEN SCANNED

=> log y

COST IN U.S. DOLLARS

SINCE FILE ENTRY

TOTAL SESSION

FULL ESTIMATED COST

0.60

0.75

STN INTERNATIONAL LOGOFF AT 15:38:18 ON 16 AUG 2000

Trying 3106016892...Open

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TERMINAL (ENTER 1, 2, 3, OR ?):2

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NEWS 6 May 22 CITED REFERENCES NOW AVAILABLE IN CAPLUS AND CA FILE

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FILE 'HOME' ENTERED AT 15:43:00 ON 16 AUG 2000

=> file reg

COST IN U.S. DOLLARS

SINCE FILE TOTAL ENTRY SESSION 0.15 0.15

FULL ESTIMATED COST

FILE 'REGISTRY' ENTERED AT 15:43:06 ON 16 AUG 2000 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
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STRUCTURE FILE UPDATES: 15 AUG 2000 HIGHEST RN 286364-83-8 DICTIONARY FILE UPDATES: 15 AUG 2000 HIGHEST RN 286364-83-8

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Please note that search-term pricing does apply when conducting SmartSELECT searches.

Structure search limits have been increased. See HELP SLIMIT for details.

=>Testing the current file.... screen

ENTER SCREEN EXPRESSION OR (END):end

=> screen 2076

L1 SCREEN CREATED

Uploading c:\stnexp4\queries\09270006.str

L2 STRUCTURE UPLOADED

=> que L2 AND L1

L3 QUE L2 AND L1

=> d 12

L2 HAS NO ANSWERS L2 STR

09/270,006

G1 O, S, P

G2 H, Cy, Hy, Ak

Structure attributes must be viewed using STN Express query preparation.

=> s 12 sss sam

SAMPLE SEARCH INITIATED 15:43:49 FILE 'REGISTRY' SAMPLE SCREEN SEARCH COMPLETED - 4609 TO ITERATE

21.7% PROCESSED 1000 ITERATIONS INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED) SEARCH TIME: 00.00.01

6 ANSWERS

FULL FILE PROJECTIONS: ONLINE **COMPLETE**

BATCH **COMPLETE**

PROJECTED ITERATIONS: 88114 TO 96246

PROJECTED ANSWERS: 238 TO 869

L4 6 SEA SSS SAM L2

=> d scan

L4 6 ANSWERS REGISTRY COPYRIGHT 2000 ACS

IN D-erythro-Hexonic acid, 2,5-anhydro-3,4-dideoxy-, benzoate (9CI)

MF C13 H14 O5

Absolute stereochemistry.

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):5

L4 6 ANSWERS REGISTRY COPYRIGHT 2000 ACS

IN 2-Furancarboxylic acid, tetrahydro-2-(1-hydroxycycloheptyl)- (9CI)
09/270,006
Page 12

MF C12 H20 O4

L4 6 ANSWERS REGISTRY COPYRIGHT 2000 ACS

IN L-Phenylalanine, (S)-tetrahydro-2-furancarboxylate (9CI)

MF C9 H11 N O2 . C5 H8 O3

CM 1

Absolute stereochemistry.

CM 2

Absolute stereochemistry. Rotation (-).

L4 6 ANSWERS REGISTRY COPYRIGHT 2000 ACS

IN [2,2':2'(3'H),2''-Terfuran]-2(3H)-carboxylic acid, octahydro-2''-[2-(octyloxy)ethyl]-, [2R*[S*(R*)]]-, compd. with 2,4,6-trinitrophenol

sodium

salt (1:1) (9CI)

MF C23 H40 O6 . C6 H3 N3 O7 . Na

CM 1

Relative stereochemistry.

CM 2

Na

L4 6 ANSWERS REGISTRY COPYRIGHT 2000 ACS

IN 2-Furanpropanoic acid, .alpha.-amino-2-carboxytetrahydro- (9CI)

MF C8 H13 N O5

L4 6 ANSWERS REGISTRY COPYRIGHT 2000 ACS

IN 2-Thiophenecarboxylic acid, tetrahydro-2-methyl-, 1,1-dioxide, (.+-.)(8CI)

MF C6 H10 O4 S

ALL ANSWERS HAVE BEEN SCANNED

 \Rightarrow log y

SINCE FILE TOTAL COST IN U.S. DOLLARS ENTRY SESSION

0.90 1.05 FULL ESTIMATED COST

STN INTERNATIONAL LOGOFF AT 15:44:39 ON 16 AUG 2000

Trying 3106016892...Open

Welcome to STN International! Enter x:x

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FILE 'HOME' ENTERED AT 15:55:40 ON 18 AUG 2000

=> file reg

COST IN U.S. DOLLARS

SINCE FILE TOTAL ENTRY SESSION 0.15 0.15

FULL ESTIMATED COST

FILE 'REGISTRY' ENTERED AT 15:55:48 ON 18 AUG 2000 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS. COPYRIGHT (C) 2000 American Chemical Society (ACS)

STRUCTURE FILE UPDATES: 17 AUG 2000 HIGHEST RN 286453-42-7 DICTIONARY FILE UPDATES: 17 AUG 2000 HIGHEST RN 286453-42-7

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Please note that search-term pricing does apply when conducting SmartSELECT searches.

Structure search limits have been increased. See HELP SLIMIT for details.

=>Testing the current file.... screen

ENTER SCREEN EXPRESSION OR (END):end

- => screen 2076
- L1 SCREEN CREATED

Uploading c:\stnexp4\queries\19270006.str

- L2 STRUCTURE UPLOADED
- => que L2 AND L1
- L3 QUE L2 AND L1
- => s 13 sss sam

09/270,006

10.5

SAMPLE SEARCH INITIATED 15:56:22 FILE 'REGISTRY' SAMPLE SCREEN SEARCH COMPLETED - 1698 TO ITERATE

58.9% PROCESSED 1000 ITERATIONS INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED) 8 ANSWERS

SEARCH TIME: 00.00.01

ONLINE **COMPLETE** FULL FILE PROJECTIONS:

COMPLETE BATCH

PROJECTED ITERATIONS:

31490 TO 36430

PROJECTED ANSWERS:

50 TO 492

8 SEA SSS SAM L2 AND L1

=> d scan

REGISTRY COPYRIGHT 2000 ACS T.4

D-erythro-Hexonic acid, 2,5-anhydro-3,4-dideoxy-, benzoate (9CI) IN

C13 H14 O5 MF

Absolute stereochemistry.

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):7

REGISTRY COPYRIGHT 2000 ACS L48 ANSWERS

IN 2-Furancarboxylic acid, tetrahydro-2-(1-hydroxycycloheptyl)- (9CI)

MF C12 H20 O4

8 ANSWERS REGISTRY COPYRIGHT 2000 ACS

[2,2':2'(3'H),2''-Terfuran]-2(3H)-carboxylic acid, octahydro-2''-[2-(octyloxy)ethyl]-, [2R*[S*(R*)]]-, compd. with 2,4,6-trinitrophenol lithium salt (1:1) (9CI)

C23 H40 O6 . C6 H3 N3 O7 . Li

CM

Relative stereochemistry.

CM 2

L48 ANSWERS REGISTRY COPYRIGHT 2000 ACS

[2,2'-Bifuran]-2(3H)-carboxylic acid, hexahydro-2'-[3-(octyloxy)propyl]-, (R*,S*)-, compd. with 2,4,6-trinitrophenol sodium salt (1:1) (9CI) C20 H36 O5 . C6 H3 N3 O7 . Na IN

MF

CM 1

Relative stereochemistry.

CM

Na

L4 8 ANSWERS REGISTRY COPYRIGHT 2000 ACS

IN [2,2':2'(3'H),2''-Terfuran]-2(3H)-carboxylic acid, octahydro-2''-[2 (octyloxy)ethyl]-, [2R*[S*(R*)]]-, compd. with 2,4,6-trinitrophenol
sodium

salt (1:1) (9CI)

MF C23 H40 O6 . C6 H3 N3 O7 . Na

CM 1

Relative stereochemistry.

CM 2

● Na

L4 8 ANSWERS REGISTRY COPYRIGHT 2000 ACS

IN [2,2':2'(3'H),2''-Terfuran]-2(3H)-carboxylic acid, octahydro-2''-[2-(octyloxy)ethyl]-, [2R*[R*(S*)]]-, compd. with 2,4,6-trinitrophenol sodium 09/270,006 Page 19

salt (1:1) (9CI) MF C23 H40 O6 . C6 H3 N3 O7 . Na

CM 1

Relative stereochemistry.

CM 2

Na

L4 8 ANSWERS REGISTRY COPYRIGHT 2000 ACS
IN 2H-Pyran-2,2,6-tricarboxylic acid, tetrahydro-, triammonium salt (9CI)
MF C8 H10 O7 . 3 H3 N

$$\begin{array}{c|c} \mathsf{HO_2C} & & \mathsf{CO_2H} \\ \hline & \mathsf{CO_2H} \\ \end{array}$$

●3 NH3

L4 8 ANSWERS REGISTRY COPYRIGHT 2000 ACS 09/270,006

IN 2H-Pyran-2-carboxylic acid, 2-(3-cyclohexen-1-yl)tetrahydro- (9CI) C12 H18 O3 MF

ALL ANSWERS HAVE BEEN SCANNED

=> s 12 sss sam

SAMPLE SEARCH INITIATED 15:58:14 FILE 'REGISTRY' SAMPLE SCREEN SEARCH COMPLETED -1698 TO ITERATE

1000 ITERATIONS 58.9% PROCESSED INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**

COMPLETE BATCH

PROJECTED ITERATIONS:

31490 TO 36430

PROJECTED ANSWERS:

50 TO 492

L5 8 SEA SSS SAM L2

=> d scan

L5 8 ANSWERS REGISTRY COPYRIGHT 2000 ACS

[2,2':2'(3'H),2''-Terfuran]-2(3H)-carboxylic acid, octahydro-2''-[2-IN (octyloxy)ethyl]-, [2R*[S*(R*)]]-, compd. with 2,4,6-trinitrophenol sodium

salt (1:1) (9CI) C23 H40 O6 . C6 H3 N3 O7 . Na MF

> CM 1

Relative stereochemistry.

09/270,006

8 ANSWERS

CM 2

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      Armour, Duncan Robert; Price, David Anthony; Stammen, Blanda Luzia
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     Osterkamp, Frank; Ziemer, Burkhard; Koert, Ulrich; Wiesner, Matthias;
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     Snider, Barry B.; Hawryluk, Natalie A.
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     Department of Chemistry, Brandeis University, Waltham, MA, 02454-9110,
USA
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     2,4-disubstituted THF rings in the molecular backbone
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     analogs over tartaric acid-modified Raney nickel
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     Nakagawa, Satoshi; Haruna, Noriko; Acosta, Delicia E.; Endo, Tadateru;
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     Faculty of Science, Himeji Institute of Technology, Hyogo, 678-1297,
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     natural cephalotaxine
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     Charbonnel, Sandra; Dhal, Robert; Dujardin, Gilles; Fournier, Florence;
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     Howard Hughes Medical Institute Department of Chemistry and Chemical
     Biology and Harvard Institute of Chemistry and Cell Biology, Harvard
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     Institute of Biological Chemistry, Academia Sinica, Taipei, 115, Taiwan
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      a]pyrimidines for inhibiting the formation of NO (nitrogen oxide) in vivo
      Yamada, Satoshi; Kinoshita, Naosumi; Yasumura, Koichi; Edamatsu, Koji;
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      Nagahama, Takao; Ishikawa, Shintaro; Yamauchi, Takeshi; Kishi, Kazumasa;
      Sugiyama, Kazuhisa
      Otsuka Pharmaceutical Co., Ltd., Japan
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     Diastereoselective .alpha.-alkoxylation of lactamide N-alkyl groups via
     intramolecular formation of oxonium ions as the key intermediate
ΑU
     Kamada, Tohru; Oku, Akira
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     Faculty of Engineering and Design, Department of Chemistry, Kyoto
     Institute of Technology, Kyoto, 606-8585, Japan J. Chem. Soc., Perkin Trans. 1 (1998), (20), 3381-3388
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      rearrangement
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      Department of Chemistry, University of Cambridge, Cambridge, CB2 1EW, UK
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     Hayes, Thomas K.; Forood, Behrouz; Kiely, John S.
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     Trega Biosciences, Inc., USA
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      PCT Int. Appl., 124 pp.
      CODEN: PIXXD2
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     Features Both Reminiscent of Natural Products and Compatible with
     Miniaturized Cell-Based Assays
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     Tan, Derek S.; Foley, Michael A.; Shair, Matthew D.; Schreiber, Stuart L.
     Department of Chemistry Chemical Biology Harvard Institute of Chemistry
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     Cambridge, MA, 02138, USA
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     Cameron, Kimberly O'Keefe; Ke, Hua Zhu; Lefker, Bruce Allen; Rosati,
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     Robert Louis; Thompson, David Duane
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     Pfizer Inc., USA
SO
     PCT Int. Appl., 276 pp.
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     Institute of Phamaceutical Research, Shanghai No. 6 Pharmaceutical
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     doxazosin, terazosin, prazosin, tiodazosin and related antihypertensive
     Zhou, Tianhao; Weeratunga, Gamini; Murthy, K. S. Keshava; Guntoori,
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     of camptothecin derivatives bearing five-membered heterocycle containing
     10-substituents
     Zhao, Rulin; Guan, Le Luo; Oreski, Bernadette; Lown, J. William
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    Dihydropyran-2-carboxylic acid, a novel bifunctional linker for the
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ΤI
     Preparation of tricyclic amide and urea compounds for inhibition of
     G-protein function and treatment of proliferative diseases
     Bishop, W. Robert; Doll, Ronald J.; Mallams, Alan K.; Njoroge, F. George;
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     Petrin, Joanne M.; Piwinski, John J.; Wolin, Ronald L.; Taveras, Arthur
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    2-(tetrahydrofuran-2'-yl)benzimidazoles
    Novelli, Federica; Tasso, Bruno; Sparatore, Fabio; Sparatore, Anna
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ΤI
     Preparation of cinchonan based chiral selectors for chiral stationary
     phases for high-performance liquid chromatography
IN
     Lindner, Wolfgang; Lammerhofer, Michael; Maier, Norbert
PA
     Lindner, Wolfgang, Austria; Lammerhofer, Michael; Maier, Norbert
SO
     PCT Int. Appl., 83 pp.
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     initiated by an inducible pyrrologuinoline quinone-dependent alcohol
     dehydrogenase
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     Zarnt, Grit; Schrader, Thomas; Andreesen, Jan R.
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ΤI
     Fluorocyclization of 2,5-tetrahydrofurandicarboxylic and
     2,3,4,5-tetrahydrofurantetracarboxylic acids with sulfur tetrafluoride
     leading to bicyclic and tricyclic fluoroethers
ΑU
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ΤI
    Process and intermediates for the synthesis of LH-RH antagonists
IN
    Funk, Kenneth W.; Lundell, Edwin O.; Miller, Robert B.; Chang, Jane L.;
    Kishore, Vimal; Napier, James J.; Staeger, Michael A.
PA
    Abbott Laboratories, USA
SO
    PCT Int. Appl., 62 pp.
    CODEN: PIXXD2
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TΙ
    Process and intermediates for the synthesis of LH-RH antagonists
    Funk, Kenneth W.; Lundell, Edwin O.; Miller, Robert B.; Chang, Jane L.;
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    Kishore, Vimal; Napier, James J.; Staeger, Michael A.
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    Abbott Laboratories, USA
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     antagonists
     Murugesan, Natesan; Barrish, Joel C.; Spergel, Steven H.
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     Bristol-Myers Squibb Company, USA
SO
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    Van Eikeren, Paul; Mc Conville, Francis X.; Lopez, Jorge
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    PCT Int. Appl., 30 pp.
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     Nakai, Sakie; Sato, Haruyo; Iwata, Shiho
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     Toray Industries, Japan
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     CODEN: JKXXAF
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     hydroxy-.alpha.-phenylacetates as potential ligands for the study of
     muscarinic receptor density by positron emission tomography
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    Health Sciences Research Division, Oak Ridge National Laboratory, Oak
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    Preparation of optically active amino acid amides as optical resolving
    agents and method for producing optically active
tetrahydrofurancarboxylic
    acid derivatives by optical resolution
    Nakai, Sakie; Sato, Haruyo; Fujino, Toshihiro
    Toray Industries, Japan
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    Peptidyl ketoamides as serine and cysteine protease inhibitors
IN
    Powers, James C.; Li, Zhaozhao; Patil, Girish S.; Chu, Der-Lun
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    Georgia Tech Research Corp, USA
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L7
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ΤI
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    Haviv, Fortuna; Fitzpatrick, Timothy D.; Swenson, Rolf E.; Nichols,
    Charles J.; Mort, Nicholas A.
    Tap Holdings Inc., USA
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    Kamioka, Chiaki; Kitagawa, Yukio
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    Shin Nippon Rika Kk, Japan
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    Synthesis and antitumor activity of camptothecin derivatives bearing
    five-membered heterocycle containing 10-substituents
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    Zhao, Rulin; Oreski, Bernadette; Lown, J. William
    Dep. Chem., Univ. Alberta, Edmonton, AB, T6G 2G2, Can.
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     G-protein function and treatment of proliferative diseases
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     Bishop, W. Robert; Doll, Ronald J.; Mallams, Alan K.; Njoroge, F. George;
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     Schering Corp., USA
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     123:228905
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     Haviv, Fortuna; Fitzpatrick, Timothy D.; Swenson, Rolf E.; Nichols,
TN
     Charles J.; Mort, Nicholas A.
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     1995:776630 CAPLUS
     123:339157
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     Kostyanovsky, R. G.; Krutius, O. N.; El'natanov, Yu. I.
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     123:305651
     Direct separation of carboxylic acid and amine enantiomers by
TΙ
     high-performance liquid chromatography on reversed-phase silica gels
     coated with chiral copper(II) complexes
     Oi, Naobumi; Kitahara, Hajimu; Aoki, Fumiko
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CS
     Sumika Chemical Analysis Service, Ltd., 3-1-135, Kasugade-naka,
     Konohana-ku, Osaka, 554, Japan
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DN
     123:313692
     Synthesis, Structure, and Pharmacological Evaluation of the Stereoisomers
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     of Furnidipine
ΑU
     Alajarin, Ramon; Vaquero, Juan J.; Alvarez-Builla, Julio; Pastor, Manuel;
     Sunkel, Carlos; Fau de Casa-Juana, Miguel; Priego, Jaime; Statkow, Peter
     R.; Sanz-Aparicio, Julia; Fonseca, Isabel
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     Departamento de Quimica Organica, Universidad de Alcala, Alcala de
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    123:70522
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ΤI
    crystal display devices
IN
    Nakazawa, Ikuo; Takiguchi, Takao; Iwaki, Takashi; Tokano, Goji; Yamada,
    Yoko; Nakamura, Shinichi
    Canon Kk, Japan
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                                      JP 1993-242160 19930903
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    Dep. Chem., Univ. Florida, Gainesville, FL, 32611, USA
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    122:81105
ΤI
    Process for preparation of N-cyanoguanidine tetrahydrofuroylamino
    derivatives via acylation of N-cyanoguanidine amines with
tetrahydrofuroyl
    mixed anhydrides
IN
    Toempe, Peter; Reiter, Jozsef; Pongo, Laszlo
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    Egis Gyogyszergyar Rt., Hung.
SO
    Hung. Teljes
    CODEN: HUXXBU
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DN
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     and esters, liquid crystal composition containing them, and liquid
crystal
     device and display apparatus
     Nakazawa, Ikuo; Nakamura, Shinichi; Takiguchi, Takao; Iwaki, Takashi;
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     Tokano, Goji; Yamada, Yoko
     Canon Kk, Japan
PA
     Jpn. Kokai Tokkyo Koho, 87 pp.
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     1995:212874 CAPLUS
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     preparation of optically active heterocyclic carboxylic acids
ΑU
     Fritz-Langhals, Elke
     Consortium Elektrochem. Ind. G.m.b.H., Munich, -D-81379, Germany
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     1994:680627 CAPLUS
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    carboxylates
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    Manero, Javier; Leupold, Ernst
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    Hoechst A.-G., Germany
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    Eur. Pat. Appl., 16 pp.
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ΤI
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    copper(II) complexes as stationary phases
ΑU
    Oi, Naobumi; Kitahara, Hajimu; Aoki, Fumiko
CS
    Sumika Chem. Anal. Serv. Ltd., Japan
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    Kuromatogurafi (1993), 14(5), 88-9
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TТ
    2-oxo-1-piperidinyl derivatives
    Ranganathan, Ramachandran S.; Arunachalam, Thangavel; Natalie, Kenneth
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J.,
    Squibb, E. R., and Sons, Inc., USA
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SO
    Eur. Pat. Appl., 28 pp.
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\mathbf{DT}
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     1994:323599 CAPLUS
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     120:323599
DN
     Oxazolidinones antibiotics containing a substituted diazine moiety
ΤI
     Hutchinson, Douglas K.; Brickner, Steven Joseph; Barbachyn, Michael
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     Robert; Gammill, Ronald B.; Patel, Mahest V.
PA
     Upjohn Co., USA
so
     PCT Int. Appl., 44 pp.
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